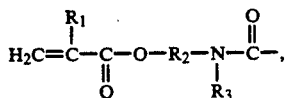


pentanediyl, neopentanediyl, hexanediyl, 2,3-dimethylbutanediyl, methylethanediyl, propanediyl and ethanediyl. Ethanediyl and propanediyl are the preferred divalent radicals.

A monovalent aliphatic C<sub>1</sub>-C<sub>4</sub>- radical (R<sub>3</sub>) in general denotes a straight-chain or branched hydrocarbon radical having 1 to 4, preferably 3 or 4, carbon atoms. The following alkyl radicals may be mentioned as examples: methyl, ethyl, propyl, isopropyl, butyl, isobutyl and tert.-butyl. tert.-Butyl and isobutyl are particularly preferred.

Z in the context of the general formula (I) denotes hydrogen (H) or an N-alkyl-N-(meth)acryloyloxyalkyl-carboxamide of the formula

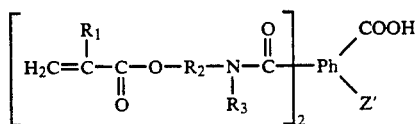


wherein R<sub>1</sub>, R<sub>2</sub> and R<sub>3</sub> have the meanings given, or also COOH, if Y also denotes COOH.

Y can also represent, in addition to COOH, an anhydride group if the two radicals containing Y are bonded to adjacent positions on an aromatic. Adjacent here means ortho-positions in the benzene or naphthalene nuclei and the α-positions (1,8- or 4,5-substitution) in the naphthalene nuclei.

Ph denotes a benzene nucleus which is trisubstituted by X and Y in the 1,2,3- or 1,2,4-position or tetrasubstituted by X, Y and Z in the 1,2,4,5-position, or a naphthalene nucleus which is trisubstituted by X and Y in the 1,2,6-, 1,4,5- or 2,3,6-position or tetrasubstituted by X, Y and Z in the 2,3,6,7- or 1,4,5,8-position.

Preferred compounds in the context of the formula (I) correspond to the formula (II)



in which

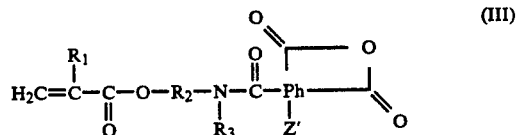
R<sub>1</sub> denotes hydrogen or methyl,  
R<sub>2</sub> denotes a divalent aliphatic radical (C<sub>2</sub>-C<sub>6</sub>),  
R<sub>3</sub> denotes a monovalent aliphatic radical (C<sub>1</sub>-C<sub>4</sub>),

Z' denotes hydrogen or COOH

and

Ph represents a tri- or tetrasubstituted benzene ring (1,2,3-/1,2,4- or 1,2,4,5-substitution) or a tri- or tetrasubstituted naphthalene ring (1,2,6-/1,4,5-/2,3,6-/1,4,5,8- or 2,3,6,7-substitution).

Another preferred group of compounds of the formula (I) corresponds to the formula (III)



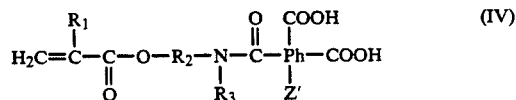
in which

R<sub>1</sub> denotes hydrogen or methyl,  
R<sub>2</sub> denotes a divalent aliphatic radical (C<sub>2</sub>-C<sub>6</sub>),  
R<sub>3</sub> denotes a monovalent aliphatic radical (C<sub>1</sub>-C<sub>4</sub>),  
Z' denotes hydrogen or COOH

and

Ph represents a benzene ring which is tri- or tetrasubstituted (1,2,3-/1,2,4- or 1,2,4,5-substitution) or a tri- or tetrasubstituted naphthalene ring (1,2,6-/1,4,5-/2,3,6-/1,4,5,8- or 2,3,6,7-substitution).

A particularly preferred group of compounds of the formula (I) corresponds to the formula (IV)



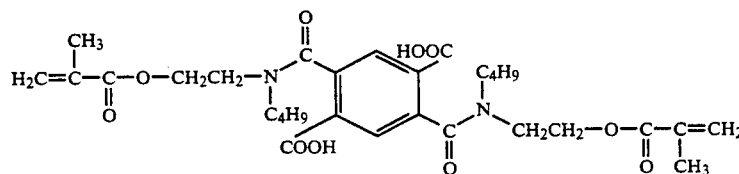
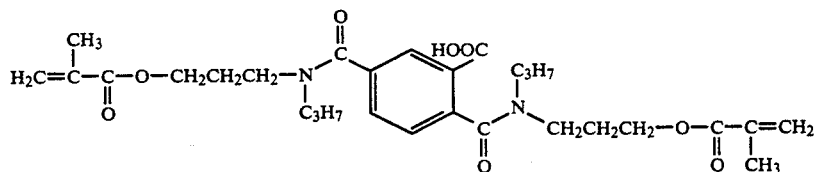
in which

R<sub>1</sub> denotes hydrogen or methyl,  
R<sub>2</sub> denotes a divalent aliphatic radical (C<sub>2</sub>-C<sub>6</sub>),  
R<sub>3</sub> denotes a monovalent aliphatic radical (C<sub>1</sub>-C<sub>4</sub>),  
Z' denotes hydrogen or COOH

and

Ph represents a tri- or tetrasubstituted benzene ring (1,2,3-/1,2,4- or 1,2,4,5-substitution) or a tri- or tetrasubstituted naphthalene ring (1,2,6-/1,4,5-/2,3,6-/1,4,5,8- or 2,3,6,7-substitution).

The following N-alkyl-N-(meth)acryloyloxyalkyl-carboxamides may be mentioned specifically as examples:



(X)